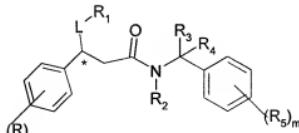


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Currently Amended) A compound of formula (I)



(I)

wherein

R is halogen, C₁₋₄ alkyl, cyano, C₁₋₄ alkoxy, trifluoromethyl or trifluoromethoxy;

R₁ is a R₁ is a 4, 5 or 6 membered heterocyclic group, wherein the 4, 5 or 6 membered

heterocyclic group may optionally be substituted by one to three substituents, which may be the same or different, selected from (CH₂)_pR₆, wherein p is zero or an integer from 1 to 4 and R₆ is selected from:

halogen,

C₁₋₄alkoxy,

C₁₋₄alkyl,

C₃₋₇cycloalkyl,

C₁₋₄ alkyl optionally substituted by halogen, cyano or C₁₋₄ alkoxy,

hydroxy,

cyano,

nitro,

trifluoromethyl,

carboxy,

NH(C₁₋₄ alkyl),

N(C₁₋₄ alkyl)₂
NH(C₃₋₇ cycloalkyl),
N(C₁₋₄ alkyl)(C₃₋₇ cycloalkyl);
NH(C₁₋₄ alkyl)OC₁₋₄ alkoxy),
OC(O)NR₇R₈,
NR₈C(O)R₇ or
C(O)NR₇R₈;

R₂ is hydrogen, or C₁₋₄ alkyl;

R₃ and R₄ independently are hydrogen, C₁₋₄ alkyl or R₃ together with R₄ and the carbon to which they are bonded is C₃₋₇ cycloalkyl;

R₅ is trifluoromethyl, S(O)_qC₁₋₄ alkyl, C₁₋₄ alkyl, C₁₋₄ alkoxy, trifluoromethoxy, halogen or cyano;

R₇ and R₈ independently are hydrogen, C₁₋₄ alkyl or C₃₋₇ cycloalkyl;

L is a single or a double bond;

n is an integer from 1 to 3;

m is zero or an integer from 1 to 3;

q is zero or an integer from 1 to 2;

provided that

a) when L is a double bond, R₁ is not an optionally substituted 5 or 6-membered heteroaryl group, in which the 5-membered heteroaryl group contains at least one heteroatom selected from oxygen, sulphur or nitrogen and the 6-membered heteroaryl group contains from 1 to 3 nitrogen atoms;

[[b]] a) the group R₁ is linked to the carbon atom shown as * via a carbon atom;
and

[[c]] b) when the heteroatom contained in the group R₁ is substituted, p is not zero; or a pharmaceutically acceptable salt thereof.

2. (Previously Presented) A compound as claimed in claim 1 wherein R is halogen or C₁₋₄ alkyl and n is an integer from 1 to 2.

3. (Previously Presented) A compound as claimed in claim 1 wherein R₅ is trifluoromethyl, methyl, methoxy, bromine, chlorine or fluorine atom and m is an integer from 1 to 2.

4. (Cancelled)

5. (Currently Amended) A compound as claimed in claim 1 wherein R is halogen or C₁₋₄ alkyl and n is an integer from 1 to 2; R₁ is piperidyl, 2-morpholinyl, 1,2,3,6-tetrahydro-4-pyridinyl, pyridyl or pyrrolidinyl and wherein R₁ is optionally substituted by one or two groups selected from halogen, C₁₋₄ alkyl or ethylC₁₋₄ alkoxy; R₂ and R₃ are independently hydrogen or methyl; R₄ is hydrogen, methyl or together with R₃ is cyclopropyl and R₅ is trifluoromethyl, methyl, methoxy, bromine, chlorine or fluorine atom and m is preferably an integer from 1 to 2.

6. (Cancelled)

7. (Currently Amended) A compound selected from N-((1*R*)-[3,5-bis(trifluoromethyl)phenyl]ethyl)-3-(4-fluorophenyl)-*N*-methyl-3-(4-piperidinyl)propionamide(diastereoisomer 1); N-((1*S*)-[3,5-bis(trifluoromethyl)phenyl]ethyl)-3-(4-fluorophenyl)-*N*-methyl-3-(4-piperidinyl)propionamide (diastereoisomer 2); N-((1*R*)-[3,5-bis(trifluoromethyl)phenyl]ethyl)-3-(4-fluorophenyl)-3-(4-fluoro-4-piperidinyl)-*N*-methylpropionamide (diastereoisomer 1); N-[(3,5-dibromophenyl)methyl]-3-(4-fluorophenyl)-3-(4-fluoro-4-piperidinyl)-*N*-methylpropionamide (enantiomer 2); N-[(3,5-bis(trifluoromethyl)phenyl)methyl]-3-(4-fluorophenyl)-3-(3-fluoro-3-piperidinyl)-*N*-methylpropionamide (diastereoisomer A); and pharmaceutically acceptable salts and solvates thereof.

8-11. (Cancelled)

12. (Previously Presented) A pharmaceutical composition comprising a compound as claimed in claim 1 in admixture with one or more pharmaceutically acceptable carriers or excipients.

13. (Cancelled)

14. (Previously Presented) A compound as claimed in claim 1 wherein R is fluorine or chlorine or methyl and n is an integer from 1 to 2.

15. (Currently Amended) A compound as claimed in claim 1 wherein R is fluorine or chlorine or methyl and n is an integer from 1 to 2; R₄ is piperidyl, 2-morpholinyl, 1,2,3,6-tetrahydro-4-pyridinyl, pyridyl or pyrrolidinyl and wherein R₁ is optionally substituted by one or two groups selected from fluorine, methyl or ethylC₁₋₄ alkoxy; R₂ and R₃ are independently hydrogen or methyl; R₄ is hydrogen, methyl or together with R₃ is cyclopropyl and R₅ is trifluoromethyl, methyl, methoxy, bromine, chlorine or fluorine atom and m is preferably an integer from 1 to 2.

16-20. (Cancelled)